

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042823\
 Data File : BP014904.D
 Acq On : 29 Apr 2023 08:11
 Operator : CG/JU
 Sample : PB152484BL
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK484

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 05/01/2023
 Supervised By : Jagrut Upadhyay 05/01/2023

Quant Time: May 01 01:12:13 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042823.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Apr 29 00:40:33 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.169	152	227798	20.000	ng/u1	0.00
20) Naphthalene-d8	11.010	136	914702	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.816	164	528035	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.569	188	1083580	20.000	ng/u1	-0.01
79) Chrysene-d12	21.651	240	930797	20.000	ng/u1	-0.01
88) Perylene-d12	24.251	264	1112194	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.464	96	37818	5.983	ng/uL	0.00
4) Pyridine-d5	3.905	84	489098	32.190	ng/u1	0.00
7) Phenol-d5	7.311	99	564319	31.652	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.487	67	371320	33.206	ng/u1	0.00
11) 2-Chlorophenol-d4	7.693	132	470275	34.837	ng/u1	0.00
15) 4-Methylphenol-d8	8.875	113	436958	32.164	ng/u1	0.00
21) Nitrobenzene-d5	9.346	128	216247	32.926	ng/u1	0.00
24) 2-Nitrophenol-d4	10.075	143	224692	37.269	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.616	165	407201	31.448	ng/u1	0.00
31) 4-Chloroaniline-d4	11.140	131	554781	29.634	ng/u1	0.00
46) Dimethylphthalate-d6	14.216	166	1297601	34.290	ng/u1	0.00
49) Acenaphthylene-d8	14.510	160	1532294	33.898	ng/u1	0.00
54) 4-Nitrophenol-d4	14.993	143	126457m	22.458	ng/u1	0.00
60) Fluorene-d10	15.804	176	1111669	33.700	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.916	200	140651	30.687	ng/u1	0.00
73) Anthracene-d10	17.669	188	1723699	34.832	ng/u1	0.00
81) Pyrene-d10	19.886	212	1935500	37.439	ng/u1	-0.01
92) Benzo(a)pyrene-d12	24.080	264	1862766	33.910	ng/u1	-0.02

Target Compounds Qvalue

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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