

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042823\
 Data File : BP014887.D
 Acq On : 28 Apr 2023 22:43
 Operator : CG/JU
 Sample : PB152473BL
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK473

Manual Integrations
 APPROVED

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

Quant Time: Apr 29 03:55:31 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	8.175	152	209897	20.000	ng/u1	0.00
20) Naphthalene-d8	11.010	136	845785	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.816	164	510420	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.575	188	1045051	20.000	ng/u1	0.00
79) Chrysene-d12	21.657	240	943102	20.000	ng/u1	0.00
88) Perylene-d12	24.257	264	1115873	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.469	96	34067	5.849	ng/uL	0.00
4) Pyridine-d5	3.905	84	374949	26.782	ng/u1	0.00
7) Phenol-d5	7.310	99	471295	28.688	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.487	67	312298	30.310	ng/u1	0.00
11) 2-Chlorophenol-d4	7.699	132	393963	31.673	ng/u1	0.00
15) 4-Methylphenol-d8	8.881	113	361462	28.876	ng/u1	0.00
21) Nitrobenzene-d5	9.351	128	181069	29.817	ng/u1	0.00
24) 2-Nitrophenol-d4	10.081	143	184821	33.154	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.622	165	322272	26.917	ng/u1	0.00
31) 4-Chloroaniline-d4	11.146	131	447245	25.837	ng/u1	0.00
46) Dimethylphthalate-d6	14.216	166	1143855	31.270	ng/u1	0.00
49) Acenaphthylene-d8	14.516	160	1364089	31.218	ng/u1	0.00
54) 4-Nitrophenol-d4	14.992	143	93617m	17.200	ng/u1	0.00
60) Fluorene-d10	15.810	176	1012627	31.757	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.916	200	98608	22.307	ng/u1	0.00
73) Anthracene-d10	17.675	188	1574562	32.991	ng/u1	0.00
81) Pyrene-d10	19.892	212	1830197	34.940	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.086	264	1826189	33.134	ng/u1	-0.01

Target Compounds Qvalue

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

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