

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP113022\
 Data File : BP012905.D
 Acq On : 01 Dec 2022 06:54
 Operator : CG/JU
 Sample : PB149320BL
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK320

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 12/01/2022
 Supervised By :mohammad ahmed 12/02/2022

Quant Time: Dec 01 07:21:02 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 29 22:53:34 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.010	152	677655	20.000	ng/u1	0.00
20) Naphthalene-d8	10.822	136	3102476	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.645	164	2152379	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.398	188	4544146	20.000	ng/u1	0.00
79) Chrysene-d12	21.480	240	3211320	20.000	ng/u1	0.00
88) Perylene-d12	23.992	264	2661261	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	84177m	5.323	ng/uL	0.00
4) Pyridine-d5	3.817	84	1171168	24.747	ng/u1	0.00
7) Phenol-d5	7.158	99	1483216	26.636	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.328	67	928475	27.059	ng/u1	0.00
11) 2-Chlorophenol-d4	7.534	132	1174961	27.691	ng/u1	0.00
15) 4-Methylphenol-d8	8.710	113	1222004	27.237	ng/u1	0.00
21) Nitrobenzene-d5	9.175	128	584472	30.780	ng/u1	0.00
24) 2-Nitrophenol-d4	9.899	143	661361	32.801	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.434	165	1251631	26.605	ng/u1	0.00
31) 4-Chloroaniline-d4	10.957	131	1755604	27.251	ng/u1	0.00
46) Dimethylphthalate-d6	14.051	166	4163625	28.104	ng/u1	0.00
49) Acenaphthylene-d8	14.339	160	4614311	26.693	ng/u1	0.00
54) 4-Nitrophenol-d4	14.828	143	566399	21.444	ng/u1	0.00
60) Fluorene-d10	15.634	176	3611679	27.836	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.745	200	475301	21.884	ng/u1	0.00
73) Anthracene-d10	17.498	188	5692147	28.457	ng/u1	0.00
81) Pyrene-d10	19.727	212	5960191	31.942	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.833	264	3764526	28.799	ng/u1	0.00

Target Compounds Qvalue

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

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