

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP120722\
 Data File : BP012974.D
 Acq On : 07 Dec 2022 17:57
 Operator : CG/JU
 Sample : PB149448BL
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK448

Quant Time: Dec 08 00:52:26 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Dec 08 00:35:07 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.981	152	532569	20.000	ng/u1	0.00
20) Naphthalene-d8	10.793	136	2304027	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.616	164	1595946	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.375	188	3708213	20.000	ng/u1	0.00
79) Chrysene-d12	21.457	240	3610164	20.000	ng/u1	0.00
88) Perylene-d12	23.945	264	4019169	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.358	96	67098	5.387	ng/uL	0.00
4) Pyridine-d5	3.787	84	926249	26.178	ng/u1	0.00
7) Phenol-d5	7.134	99	1119104	25.798	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.304	67	730061	27.899	ng/u1	0.00
11) 2-Chlorophenol-d4	7.504	132	949076	27.793	ng/u1	0.00
15) 4-Methylphenol-d8	8.681	113	928010	26.109	ng/u1	0.00
21) Nitrobenzene-d5	9.146	128	473486	27.970	ng/u1	0.00
24) 2-Nitrophenol-d4	9.869	143	537814	28.220	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.410	165	955092	25.801	ng/u1	0.00
31) 4-Chloroaniline-d4	10.928	131	1422123	27.213	ng/u1	0.00
46) Dimethylphthalate-d6	14.022	166	3588433	29.256	ng/u1	0.00
49) Acenaphthylene-d8	14.310	160	3777835	28.030	ng/u1	0.00
54) 4-Nitrophenol-d4	14.804	143	516195	23.396	ng/u1	0.00
60) Fluorene-d10	15.610	176	3007954	28.988	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.722	200	560297	23.480	ng/u1	0.00
73) Anthracene-d10	17.469	188	4949326	29.584	ng/u1	0.00
81) Pyrene-d10	19.698	212	5987278	28.893	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.786	264	5891110	29.407	ng/u1	0.00

Target Compounds Qvalue

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

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